

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061518\
 Data File : BF106492.D
 Acq On : 15 Jun 2018 15:28
 Operator : JU/SJ
 Sample : J3471-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 #2

Manual Integrations
 APPROVED

Sohil
 6/18/2018 10:38:25 AM

Quant Time: Jun 16 00:04:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 15 11:11:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	113186	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	410916	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	176798	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	289974	20.00	ng	0.00
75) Chrysene-d12	14.15	240	303105	20.00	ng	0.00
86) Perylene-d12	15.71	264	241878	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	615728	92.07	ng	0.00
7) Phenol-d6	6.61	99	740610	87.66	ng	0.00
23) Nitrobenzene-d5	7.52	82	486648	69.67	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	159768	93.92	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	757494	72.53	ng	0.00
78) Terphenyl-d14	13.08	244	649752	53.24	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	78842	6.242	ng	99
70) Phenanthrene	11.52	178	47907	3.373	ng	98
74) Fluoranthene	12.71	202	121542	8.025	ng	96
77) Pyrene	12.94	202	120397	6.345	ng	99
80) Benzo(a)anthracene	14.14	228	67850	3.762	ng	95
82) Chrysene	14.18	228	69240	4.204	ng	97
85) Indeno(1,2,3-cd)pyrene	17.27	276	27252	2.073	ng	# 88
87) Benzo(b)fluoranthene	15.25	252	78036m	5.337	ng	
89) Benzo(a)pyrene	15.64	252	51128	3.888	ng	96
91) Benzo(g,h,i)perylene	17.75	276	29691	2.787	ng	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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